AMENDMENTS TO THE CLAIMS

This listing of claims will replace all prior versions, and listings, of claims in the application:

Claims 1-36 (cancelled)

37. (previously presented) A compound represented by formula (1) according to claim 56, in which R5 is a phenyl group substituted at least in position 3.

38. (previously presented)

A compound represented by formula (1) according to claim 56, wherein R5 is a phenyl group substituted at least in position 3 by a substituent selected from the group consisting of: CHO, CN, CONH2, NO2, CF3, NH2, halogen, (C1-C6) alkyl, a phenyl group optionally substituted by acetyl, halogen, CONH2 or CN, a prop-1-ynyl group optionally substituted, by benzyloxy or butyl carbamate, a hex-1-ynyl group optionally substituted by CN or NH2, a pentyl group optionally substituted by CONH2, a hexyl group, a piperidinyl group optionally substituted by prop-1-ynyl, benzylaminomethyl, acetamide, aminomethyl, NH2CS-, 4-phenyl-1,3-thiazo-2-yl, -CONHBenzyl, -COOEthyl, -CONHiPropyl, -CONH-(CH2)n-CONH2 (wherein n represents a whole number from 1 to 6), -CONR'R" wherein R' and R", which are the same or different, represent C1 to C6 alkyl or hydrogen, -(4-benzylpyperazin-1-yl)carbonyl, -CONH-(CH2)n-phenyl (wherein n represents a whole number from 1 to 6), imidazolyl, or a piperazinyl group optionally substituted by a phenyl group.

39. (previously presented) A compound represented by formula (1) according to claim 56, in which R5 is a phenyl group substituted in positions 3 and 4 by a hydrocarbon chain optionally containing at least one heteroatom, said chain forming a ring with the phenyl group to which it is attached.

40. (previously presented) A compound represented by formula (1) according to claim 56, in which R5 is 3-pyridyl, 4-isoquinolyl or a_piperazinyl group optionally substituted in position 4 by an aryl group.

41. (previously presented) A compound represented by formula (1) according to claim 56, in which Z represents a sulfur atom or -NR2, wherein R2 and R1 form an imidazole.

42. (cancelled)

43. (previously presented) A compound represented by formula (1) according to claim 56, in which R1 represents hydrogen (C_1 - C_6) alkyl, (C_6 - C_{18}) aryl, (C_6 - C_{18}) aryl(C_1 - C_4) alkyl, or a (C_1 - C_{12}) alkyl(C_6 ,- C_{18}) aryl group, said aryl group being optionally substituted.

44. (previously presented) A compound represented by formula (1) according to claim 56, in which R5 is a phenyl group substituted by a group selected from:

- (a) one or more OR' groups,
- (b) a COR' group, or
- (c) a CONR'R" group
- (d) a CN group,
- (e) a trifluoromethyl group,
- (f) an alkyl group or alkynyl group,
- (g) an aryl or heterocycle group optionally substituted by one or more groups selected from groups (a)-(g).

- 45. (previously presented) A compound represented by formula (1) according to claim 56, in which at least one of R_7 and R_8 represents OR_{10} wherein R10 represents a (C_1-C_6) alkyl or (C_3-C_5) cycloalkyl group.
- 46. (previously presented) A compound represented by formula (1) according to claim 56, in which at least one group R_7 , or R_8 represents OR_{10} wherein R_{10} represents a (C_1 - C_6) alkyl or (C_3 - C_6) cycloalkyl group, and at least one of R_7 and R_8 represents a methoxy group.
- 47. (currently Amended) A compound according to claim 56, wherein said compound is selected from the group consisting of:
- 3-(7,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile, **3a** 7,8-dimethoxy-[5-(3-trifluoromethyl)phenyl]-1,3-dihydro-2H-1,4-benzodiazepin-2-one, **3d**
- 3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-benzonitrile, **4a**
- 3-[1-(4-chlorobenzyl)-7,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl]-benzonitrile, ${f 4c}$
- 3-[1-(3,4-chlorobenzyl)-7,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl]-benzonitrile, ${\bf 4d}$
- 3-[7,8-dimethoxy-1-(4-methoxybenzyl)-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl]-benzonitrile, $\bf 4e$
- 3-[1-(3-chlorobenzyl)-7,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl]-benzonitrile, $\bf 4f$
- 3-(7,8-dimethoxy-2-oxo-1-[3-(trfluoromethyl)benzyl]-2,3-dihydro-1H-1,4-benzodiazepin-5-yl]-benzonitrile, **4g**
- 3-[1-(2-chiorobenzyl)-7,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl]-benzonitrile, ${\bf 4h}$
- 3-{7,8-dimethoxy-2-oxo-1-[4-(trifluoromethyl)benzyl]-2,3-dihydro-1H-1,4-benzodiazepin-5-yl]-benzonitrile, $\bf 4i$
- 3-[7,8-dimethoxy-2-oxo-1-(2-phenylethyl)-2,3-dihydro-1H-1,4-benzodiazepin-5-yl]-

5f3-(1-ethyl-7,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzamide, **5g**3-(1-benzyl-7,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzamide, **5h**

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Serial No. 10/533,157
Art Unit 1624
ethyl{5-[3-(aminocarbonyl)phenyl]-7,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-
benzodiazepin-1-yl}acetate, 5i
3-(7,8-dimethoxy-1,3-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yi)
benzamide, 5i
3-[3-(3,4-dichiorobenzyl)-7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-
benzodiazepin-5-yl]benzamide, 5k
3 (8-methoxy-1-methyl-2 oxo-2,3 dihydro-1H-1,4-benzodiazepin-5-yl)benzamide, 5l
3-(7,8-dimethoxy-1-methyl-2-oxo-9-phenyl-2,3-dihydro-1H-1,4-benzodiazepin-
5yi)benzamide, 5m
3 (6,8-dimethoxy 2 oxo 2,3-dihydro 1H,4-benzodiazepin 5 yl)benzamide, 5n
3 (6,8 dimethoxy 1 methyl 2 oxo 2,3 dihydro lH 1,4 benzodiazepin 5 yl)benzamide,
50
tert-butyl-3-[3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-
yl)phenyl]propynylcarbamate. 6a
7,8-dimethoxy-5-(3'-hex-1-ynylphenyl)-1-N-methyl-1,3-dihydro-2H-1,4-benzodiazepin-
2-one, 6b
7,8-dimethoxy-l-methyl-5-[3-(3-piperidin-1-ylprop-1-ynyl)phenyl]-1,3-dihydro-2H-1,4-
benzodiazepin-2-one, 6c
6-[3-(7,8-dimethoxy-l-methyl-2-oxo-2.3-dihydro-lH-1,4-benzodiazepin-5-
yl)phenyl]hex-5-ynenitrile, 6d
7,8-dimethoxy-5-(3'-hexylphenyl)-1-N-methyl-1,3-dihydro-2H1,4 benzodiazepin-2-one,
6e
5-[3-(3-aminopropyl)phenyl-7,8-dimethoxy-1-methyl-1,3-dihydro-2H-1,4-
benzodiazepin-2-one trifluoroacetate, 6h
6-[3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodlazepin-5-
yl)phenyl]hexanamide, 6i
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5-{3-[3-(benzyloxy)prop-l-ynyl]phenyl}-1-ethyl-7,8-dimethoxy-1,3-dihydro-2H-1,4-benzodiazepin-2-one, $\bf 6k$

5-(4'-chloro-1,1'-biphenyl-3-yl)-7,8-dimethoxy-l-methyl-1,3-dihydro-2H-1,4-

benzodiazepin-2-one, 6j

- 3'-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-1,1'-biphenyl-3-carbonitrile, **61**
- 3'-(7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-1,1'-biphenyl-4-carbonitrile, $\bf 6m$
- 3'-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-1,1'-biphenyl-4-carboxamide, $\bf 6n$
- 3'-(7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-1,1'-biphenyl-3-carboxamide, $\bf 6o$
- 3-[3-(3,4-dichlorobenzyl)-7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl]benzonitrile, $7\mathbf{b}$
- 7,8-dimethoxy-1,3-dimethyl-5-(3-trifluoromethylphenyl)-1,3-dihydro-2H-1,4-benzodiazepin-2-one, 7c
- 3-(7,8-dimethoxy-1,3-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile, $7\mathbf{d}$
- $5-[3-(aminomethyl)phenyl]-7,8-dimethoxy-1-methyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one, {\bf 8a}$
- N-[3-(7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1 H-1,4-benzodiazepin-5-yl)benzyl]acetamide, ${\bf 8b}$
- 3-(7,8-dlmethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)thiobenzamide, $\bf 9a$
- 7,8-dimethoxy-1-methyl-5-[3-(4-phenyl-1,3-thiazol-2-yl)phenyl]-1,3-dihydro-2H-1,4-benzodiazepin-2-one, $\bf 9b$
- $5\hbox{-}(3\hbox{-}cyanophenyl)\hbox{-}7,8\hbox{-}dimethoxy\hbox{-}1,3\hbox{-}dihydro\hbox{-}2H\hbox{-}1,4\hbox{-}benzodiazepin\hbox{-}2\hbox{-}thione,} \textbf{10d}$
- 3 (8,9 dimethoxy 4H-imidazo[1,2-a][1,4]benzodiazepin-5-yl)benzonitrile, 11a
- 3 (8,9 dimethoxy-4H-imidazo[1,2-a)[1,4]benzodiazepin-6-yl)benzamide, 11b
- 3-(7,8-dimethoxy-2-methylamino-1,3-dihydro-3H-1,4-benzodiazepin-5-yl)benzonitrile, **12a**
- 7,8-dimethoxy-l-methyl-5-(3-pyridyl)-1,3-dihydro-1,4-benzodiazepin-2-one, **17b** 7,8-dimethoxy-l-methyl-5-(3-nitrophenyl)-1,3-dihydro-1,4-benzodiazepin-2-one, **17c** 5-(7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-benzonitrile, **17d**

- $5-(3-acetylphenyl)-7,8-dimethoxy-l-methyl-1,3-dihydro-1,4-benzodiazepin-2-one, \ensuremath{\mathbf{17e}}\xspace$ $5-(4-isoquinolinyl)-7,8-dimethoxy-l-methyl-1,3-dihydro-1,4-benzodiazepin-2-one, \ensuremath{\mathbf{17f}}\xspace$ 7,8-dimethoxy-5-(3-hydroxymethylphenyl)-1-methyl-3-propyl-1,3-dihydro-2H-1,4 benzodiazepin-2-one, $\ensuremath{\mathbf{17h}}\xspace$ benzodiazepin-2-one, $\ensuremath{\mathbf{17h}}\xspace$
- 5-(3-aminophenyl)-7,8-dimethoxy-1-methyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one, 17i
- 5-(3,4-dichlorophenyl)-7, 8-dimethoxy-1-methyl-1,3-dihydro-1,4-benzodiazepin-2-one, ${f 17J}$
- 7,8-dimethoxy-l-methyl-5-(3-methylphenyl)-1,3-dihydro-1,4-benzodiazepin-2-one, **17k**. 5-(3-formyiphenyly-7,8-dimethoxy-l-methyl-1,3-dihydro-1,4-benzodiazepin-2-one, **171** 5-[3-(benzylaminomethyl)phenyl]-7,8-dimethoxy-1-methyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one hydrochloride, **17m**
- $N-[3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)\\ phenyl] acetamide, {\bf 17n}$
- 7,8-dimethoxy-1-methyl-5-(3,4-methylenedioxyphenyl)-1,3-dihydro-2H-1,4-benzodiazepin-2-one, **170**
- 3-(7-hydroxy-8-methoxy-2-oxo-2,3-dihydro-111,4-benzodiazepin-5-yl)benzonitrile, **22b** 3-(6-bromo-7-hydroxy-8-methoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile, **23b**
- 3-(9-bromo-8-hydroxy-7-methoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile, **23d**
- 3-(6-bromo-7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile, **24b**
- 3-(7,8-dimethoxy-1-methyl-2-oxo-6-phenyl-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile, **25b**
- 3-(7,8-dimethoxy-1-methyl-2-oxo-9-phenyl'-2,3-dihydro-1H-1,4rbenzodiazepin-5-yl)benzonitrile,**25a**
- $tert\text{-butyl-3-[5-(cyanophenyl)-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-9-yl)phenyl] prop-2-ynylcarbamate, \textbf{25c}$
- $methyl(2E)-3-[5-(cyanophenyl)-7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-9-yl)phenyl] acrylate, \begin{center} \bf 25d \end{center}$

- tert-butyl-3-[5-(cyanophenyl)-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1 H-1,4-benzodiazepin-6-yl)phenyl]prop-2-ynylcarbamate, **25e**
- [9-(3-aminoethynyl)-7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-H-1,4-benzodiazepin-5-yl]benzonitrile, ${\bf 25f}$
- [6-(3-aminoethynyl)-7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl]benzonitrile, **25g**
- 3 (8 methoxy 2 oxo 2,3 dihydro 1H-1,4 benzodiazepin 5 yl)benzonitrile, 28a
- 3 (6-methoxy 2 oxo 2,3 dihydro 1H 1,4-benzodiazepin 5-yl)benzonitrile, 28b
- 3 (7-methoxy 2 oxo 2,3 dihydro 1H-1,4 benzodiazepin 5 yi)benzonitrile, 28e
- 6 methoxy-5-phenyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one, 28d
- 7 methoxy 5 phenyl 1,3 dihydro 2H 1,4 benzodiazepin 2 one, 28e
- 9-bromo-7,8-dimethoxy-5-phenyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one, 28f
- 3 (6,8-dimethoxy 2 oxo 2,3-dihydro 1H-1,4-benzodiazepin-5-yl)benzonitrile, 28g
- $3\hbox{-}(7\hbox{-}bromo-6, 8\hbox{-}dimethoxy-2\hbox{-}oxo-2, 3\hbox{-}dihydro-1H-1, 4\hbox{-}benzodiazepin-5-yl)} benzonitrile,$

28h

- 3 (6,8 methoxy l methyl 2 oxo 2,3 dihydro 1H 1,4 benzodiazepin 5 yl)benzonitrile, 29a
- 3-(6,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile, **29b**
- 3-(7-bromo-6,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile, $\bf 29c$
- 3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)methyl benzoate, **34a**
- 3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzoic acid, **35a**
- 3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1 H-1,4-benzodiazepin-5-yl)N-isopropylbenzamide, **36a**
- N-benzyl-3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl) benzamide, ${\bf 36b}$
- N-(6-amino-6-oxohexyl)-3-(7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzamide, **36c**

- 3-(7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-N,N-dimethylbenzamide ${\bf 36d}$
- $5-(3-[(4-benzylpiperazin-1-yl)carbonyl]phenyl)_{7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-2-one, $36e$ and$
- 3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro- 1 H-1,4-benzodiazepin-5-yl)-N-(3-phenylpropyl)benzamide, **36f**.
- 48. (previously presented) The compound according to claim 36, wherein said compound is selected from the group consisting of:
- 3-(1-benzyl-7,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile, **4m**
- 7,8-dimethoxy-1-methyl-[5-(3-trifluoromethyl)phenyl]-1,3-dihydro-2H-1,4-benzodiazepin-2-one, $\bf 4p$
- 3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzamide, **5a**
- 3-(6-bromo-7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzamide, ${f 5b}$
- tert-butyl-3-[3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)phenyl]propynylcarbamate, $\bf 6a$
- 7,8-dimethoxy-5-(3'-hex-1-ynylphenyl)-1-N-methyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one, ${\bf 6b}$
- 6-[3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)phenyl]hex-5-ynenitrile, **6d**
- 7,8-dimethoxy-5-(3'-hexylphenyl)-1-N-methyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one, $\mathbf{6e}$
- 5-(4'-chloro-1,1'-biphenyl-3-yl)-7,8-dimethoxy-l-methyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one, $\bf{6j}$
- 3'-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-1,1'-biphenyl-4-carbonitrile, ${\bf 6m}$
- 3'-(7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1 H-1,4-benzodiazepin-5-yl)-1,1'-biphenyl-4-carboxamide, $\bf 6n$

3-(3,4-dichlorobenzyl)-1-ethyl-7,8-dimethoxy-5-phenyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one, 7a

7,8-dimethoxy-1-methyl-5-[3-(4-phenyl-1, 3-thiazol-2-yl)phenyl]-1,3-dihydro-2H-1,4-benzodiazepin-2-one, $\bf 9b$ and

7,8-dimethoxy-l-methyl-5-(3-pyridyl)-1,3-dihydro-1,4-benzodiazepin-2-one, 17b.

49. (cancelled)

Claims 50-54 (cancelled)

55. (currently amended) A compound represented by formula (1) according to claim 56 [[39]], wherein R5 is a phenyl group substituted in positions 3 and 4 by a methylenedioxy chain that forms a ring with said phenyl group.

56. (currently amended) A compound represented by formula (1)

$$\begin{array}{c|c}
R_8 & R_1 \\
\hline
R_7 & R_3
\end{array}$$

$$\begin{array}{c|c}
R_8 & R_5
\end{array}$$

$$\begin{array}{c|c}
R_1 & Z \\
R_3 & R_3
\end{array}$$

$$\begin{array}{c|c}
R_1 & Z \\
R_2 & R_3
\end{array}$$

$$\begin{array}{c|c}
R_1 & Z \\
R_2 & R_3
\end{array}$$

$$\begin{array}{c|c}
R_3 & R_3
\end{array}$$

$$\begin{array}{c|c}
R_1 & R_3
\end{array}$$

in which:

Z represents oxygen, sulfur or -NR₂;

 R_1 is selected from the group consisting of hydrogen, (C_1-C_6) alkyl, (C_6-C_{18}) aryl,- (C_1-C_6) alkyl (C_6-C_{18}) aryl and (C_6-C_{18}) aryl (C_1-C_4) alkyl;

 R_2 is selected from the group consisting of hydrogen, (C_1-C_6) alkyl, (C_6-C_{18}) aryl, (C_1-C_6) alkyl (C_6-C_{18}) aryl and (C_6-C_{18}) aryl (C_1-C_4) aryl, and wherein Z is -NR2, R_1 and R_2 taken together may form a linear- or branched-hydrocarbon chain having from 2 to 6 carbon

atoms, optionally containing one or more double bonds and optionally containing oxygen, sulfur or nitrogen;

 R_3 and R_3 ', are independently selected from the group consisting of hydrogen, - (C_1C_{12}) alkyl, $(C_3$ - $C_6)$ cycloalkyl, $(C_6$ - $C_{18})$ aryl, $(C_6$ - $C_{18})$ aryl, $(C_1$ - $C_4)$ alkyl, $(C_1$ - $C_{12})$ alkyl, $(C_6$ - $C_{18})$ aryl, $(C_5$ - $C_{18})$ heterocycle, aromatic or not, containing 1 to 3 heteroatoms, - NO_2 , CF_3 , CN, NR'R'', SR', OR', COOR' and CONR'R'', wherein R' and R'' are independently selected from the group consisting of hydrogen, $(C_1$ - $C_6)$ alkyl, $(C_3$ - $C_6)$ cycloalkyl, $(C_6$ - $C_{12})$ aryl and a $(C_5$ - $C_{12})$ heterocycle, aromatic or not, containing 1 to 3 heteroatoms;

R₅ represents a phenyl group substituted at least in position 3; a naphthyl group; a (C₅-C₁₈) heterocycle, aromatic or not, containing 1 to 3 heteroatoms, selected from the group consisting of pyridyl, isoquinolyl, quinolyl and piperazinyl, provided that, when R_5 is a naphthyl group substituted in position 6, said naphthyl group is not attached to the rest of the molecule at position 2, or when R₅ is a pyridyl group, said pyridyl group is not attached to the rest of the molecule at position 4, or when R₅ is a tetrahydro-1,2,3,4isoquinolyl group, said group is not attached to the rest of the molecule at position 2, or when R₅ represents a phenyl group substituted at least in position 3, said substituent is selected from the group consisting of: alkyl, halogenoalkyl, cycloalkyl, alkenyl, alkynyl, aralkyl, aryl, (C5-C18) heterocycle, aromatic or not, hydroxy, = O, NO2, NH2, CN, COR', COOR', (C1-C6)alkoxy, (di)(C1-C6)alkylamino, NHCOR', CONR'R" group, in which R' and R" are as defined hereinabove, CHO, CONH2 and phenyl optionally substituted by an acetyl group, halogen, CONH2 or CN; prop-1-ynyl optionally substituted by a benzyloxy or tert-butyl carbamate; hex-1-ynyl optionally substituted by CN or NH2; pentyl optionally substituted by CONH2; hexyl; piperidinyl optionally substituted by a prop-1ynyl, benzylaminomethyl, acetamide, aminomethyl, NH2CS-, 4-phenyl-1,3-thiazol-2-yl, -CONHBenzyl, -COOEthyl, -CONHiPropyl, -CONH-(CH2)n-CONH2 (wherein n represents a whole number from 1 to 6), a -CONR'R" group, wherein R' and R" are independently C1-C6 alkyl or hydrogen, -(4-benzylpyperazin-1-yl)carbonyl, -CONH-(CH2)n-phenyl (wherein n represents a whole number from 1 to 6), imidazolyl, piperazinyl optionally substituted by a phenyl group, or R5 is substituted in positions 3

and 4 by halogens or by a hydrocarbon chain, optionally containing a heteroatom, thereby forming a ring with the phenyl;

 R_7 and R_8 are independently selected from the group consisting of hydrogen, halogen and a OR_{10} , group in which R_{10} represents hydrogen, (C_1-C_6) alkyl, (C_3-C_6) cycloalkyl, (C_6-C_{12}) aryl, or a (C_5-C_{12}) heterocycle, aromatic or not, containing 1 to 3 heteroatoms, wherein at least one R_7 and R_8 represents a OR_{10} group as defined hereinabove;

 R_6 and R_9 are independently selected from the group consisting of hydrogen, halogen, alkyl, cycloalkyl, alkenyl, alkynyl, aryl, aralkyl, heterocycle, aromatic or not, and an OR_{10} group, wherein R10 is as defined hereinabove;

further wherein the alkyl, cycloalkyl, alkenyl, alkynyl, aralkyl, aryl and heterocycle, aromatic or not, groups, or the hydrocarbon chain defined hereinabove are optionally substituted by one or more substituents independently selected from the group consisting of halogen, alkyl, halogenoalkyl, cycloalkyl, alkenyl, alkynyl, aralkyl, aryl, heterocycle, aromatic or not, hydroxy, =0, NO₂, NH₂, CN, COR', COOR', (C_1 - C_6)alkoxy, (di)(C_1 - C_6)alkylamino, NHCOR' and CONR'R" groups, in which R' and R" are as defined hereinabove,

or [[and]] pharmaceutically acceptable salts thereof,

with the exception of compounds represented by formula (1) in which:

R1 is alkyl, R3, R3', R6 and R9 are hydrogen, and R5 is a phenyl group substituted at least in position 3 by a methoxy group;

R1 is alkyl or hydrogen, R3, R3', R6 and R9 are hydrogen, R5 is a phenyl group substituted only in position 3 by chlorine or bromine;

R1 is alkyl, R3, R3', R6 and R9 are hydrogen, R5 is a phenyl group substituted at least in position 3 by a CH2OH group;

R1, R3, R3', R6 and R9 are hydrogen, R5 is a phenyl group substituted only in position 3 by CF3,

R1 is alkyl, R3, R3', R6 and R9 are hydrogen, R5 is a phenyl group substituted in positions 3 and 5 by CF3,

R1 is alkyl, R3, R3', R6 and R9 are hydrogen, R7 and R8 are methoxy, R5 is a phenyl group substituted in position 3 by a phenyl group, and

R1 is alkyl, R3, R3', R6 and R9 are hydrogen, R7 and R8 are methoxy, R5 is a phenyl group substituted in position 3 by a phenylethynyl group.

57. (currently amended) The compound of claim 44 wherein R5 is <u>phenyl substituted</u> with OR' wherein R' is an alkyl or alkynyl.

58. (previously amended) The compound of claim 46 wherein said compound is 7,8-dimethoxy-5-(3'-hexylphenyl)-1-N-methyl-1,3-dihydro-2H1,4 benzodiazepin-2-one.